

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012787.D
 Acq On : 27 Nov 2022 02:41
 Operator : CG/JU
 Sample : N5661-02
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C9803

Quant Time: Nov 27 23:02:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	478085	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	2171171	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	1488983	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	3322848	20.000	ng/u1	0.00
79) Chrysene-d12	21.492	240	3341784	20.000	ng/u1	0.00
88) Perylene-d12	24.010	264	3360668	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	55665	4.989	ng/uL	0.00
4) Pyridine-d5	3.823	84	223000	6.679	ng/u1	0.00
7) Phenol-d5	7.164	99	259002	6.593	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	791566	32.699	ng/u1	0.00
11) 2-Chlorophenol-d4	7.546	132	621074	20.747	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	433955	13.710	ng/u1	0.00
21) Nitrobenzene-d5	9.181	128	481929	36.266	ng/u1	0.00
24) 2-Nitrophenol-d4	9.910	143	431856	30.606	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	773521	23.495	ng/u1	0.00
31) 4-Chloroaniline-d4	10.963	131	1130494	25.075	ng/u1	0.00
46) Dimethylphthalate-d6	14.057	166	3516919	34.315	ng/u1	0.00
49) Acenaphthylene-d8	14.345	160	3957676	33.095	ng/u1	0.00
54) 4-Nitrophenol-d4	14.828	143	80603	4.411	ng/u1	0.00
60) Fluorene-d10	15.645	176	3290090	36.656	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	376608	23.714	ng/u1	0.00
73) Anthracene-d10	17.504	188	5699785	38.968	ng/u1	0.00
81) Pyrene-d10	19.733	212	6777636	34.905	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.845	264	6515930	39.473	ng/u1	0.00

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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